

conducting SmartSELECT searches.

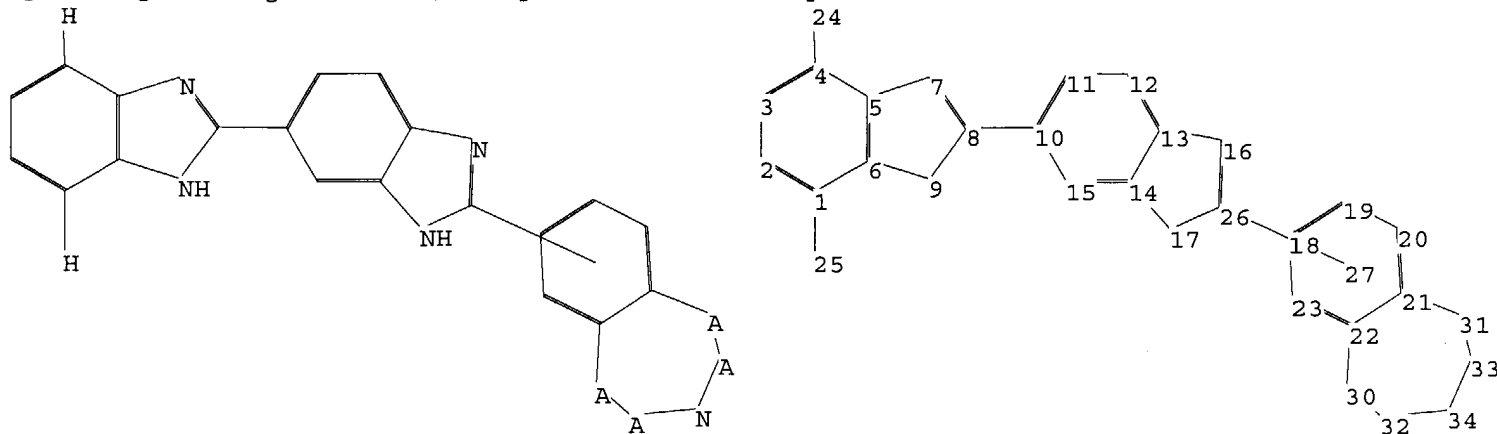
Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer

to the file summary sheet on the web at:

<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>  
Uploading C:\Program Files\Stnexp\Queries\10690800j.str



chain nodes :

24 25

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 26 30 31  
32 33 34

chain bonds :

1-25 4-24 8-10

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 10-11 10-15 11-12 12-13 13-14 13-16  
14-15 14-17 16-26 17-26 18-23 18-19 19-20 20-21 21-22 21-31 22-23 22-30 30-32  
31-33 32-34 33-34

exact/norm bonds :

5-7 6-9 7-8 8-9 13-16 14-17 16-26 17-26 21-31 22-30 30-32 31-33 32-34 33-34

exact bonds :

1-25 4-24 8-10

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 10-11 10-15 11-12 12-13 13-14 14-15 18-23 18-19 19-20  
20-21 21-22 22-23

G1

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom  
12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom  
22:Atom 23:Atom 24:CLASS 25:CLASS 26:Atom 27:CLASS 30:Atom 31:Atom 32:Atom 33:Atom  
34:CLASS

L1 STRUCTURE UPLOADED

=> s l1

SAMPLE SEARCH INITIATED 14:04:44 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 54 TO ITERATE

100.0% PROCESSED 54 ITERATIONS

SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 640 TO 1520  
PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s l1 full

FULL SEARCH INITIATED 14:04:47 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 1067 TO ITERATE

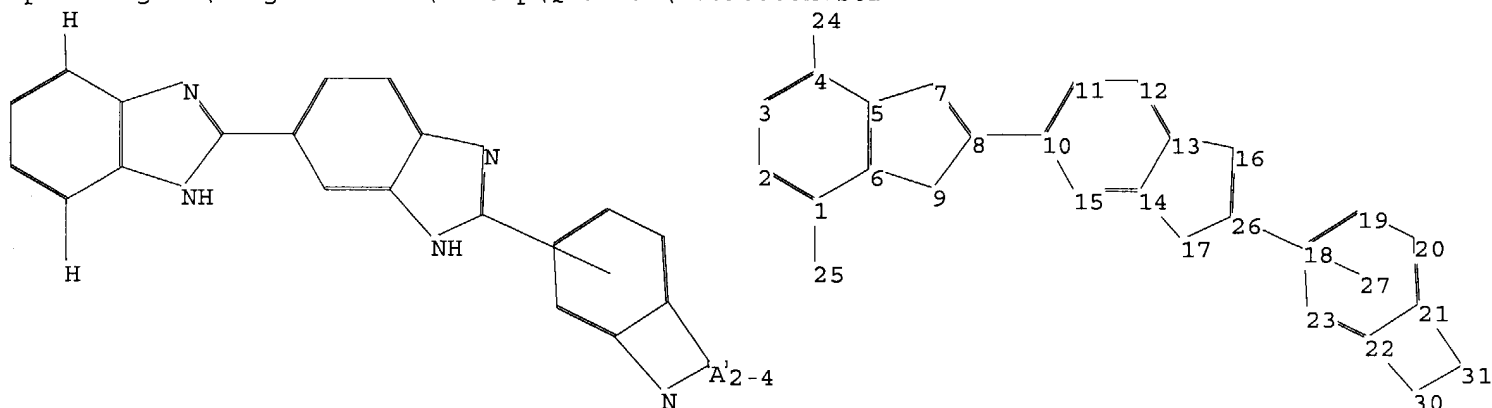
100.0% PROCESSED 1067 ITERATIONS  
SEARCH TIME: 00.00.01

0 ANSWERS

L3 0 SEA SSS FUL L1

=>

Uploading C:\Program Files\Stnexp\Queries\10690800h.str



chain nodes :

24 25

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 26 30 31

chain bonds :

1-25 4-24 8-10

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 10-11 10-15 11-12 12-13 13-14 13-16  
14-15 14-17 16-26 17-26 18-23 18-19 19-20 20-21 21-22 21-31 22-23 22-30 30-31

exact/norm bonds :

5-7 6-9 7-8 8-9 13-16 14-17 16-26 17-26 21-31 22-30 30-31

exact bonds :

1-25 4-24 8-10

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 10-11 10-15 11-12 12-13 13-14 14-15 18-23 18-19 19-20  
20-21 21-22 22-23

G1

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom  
12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom  
22:Atom 23:Atom 24:CLASS 25:CLASS 26:Atom 27:CLASS 30:Atom 31:Atom

L4 STRUCTURE UPLOADED

=> s l4

SAMPLE SEARCH INITIATED 14:05:17 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 14 TO ITERATE

100.0% PROCESSED 14 ITERATIONS  
SEARCH TIME: 00.00.01

4 ANSWERS

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 56 TO 504  
PROJECTED ANSWERS: 4 TO 200

L5 4 SEA SSS SAM L4

=> s l4 full

FULL SEARCH INITIATED 14:05:20 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 364 TO ITERATE

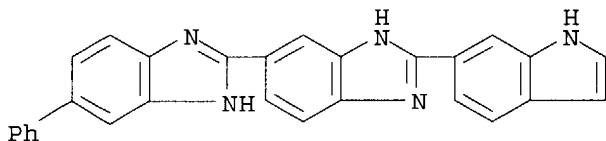
100.0% PROCESSED 364 ITERATIONS  
SEARCH TIME: 00.00.01

75 ANSWERS

L6 75 SEA SSS FUL L4

=> d scan

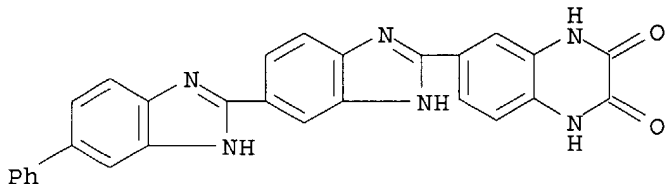
L6 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN 2,5'-Bi-1H-benzimidazole, 2'-(1H-indol-6-yl)-5-phenyl- (9CI)  
MF C28 H19 N5



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L6 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN 2,3-Quinoxalinedione, 1,4-dihydro-6-(5-phenyl[2,5'-bi-1H-benzimidazol]-2'-yl)- (9CI)  
MF C28 H18 N6 O2

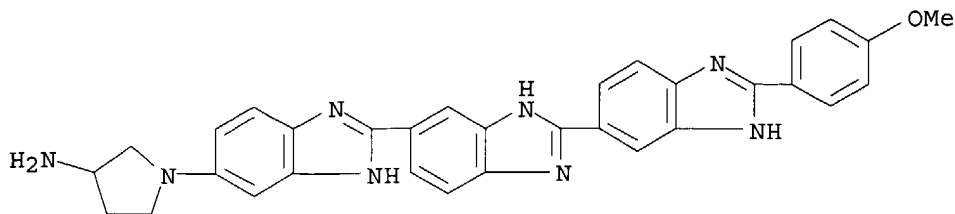


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L6 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN Guanosine, 2'-deoxycytidylyl-(5'→3')-2'-deoxyguanylyl-(5'→3')-2'-deoxycytidylyl-(5'→3')-2'-deoxyadenylyl-(5'→3')-2'-deoxyadenylyl-(5'→3')-2'-deoxyadenylyl-(5'→3')-thymidylyl-(5'→3')-thymidylyl-(5'→3')-2'-deoxyguanylyl-(5'→3')-2'-deoxycytidylyl-(5'→3')-2'-deoxy-, double-stranded complementary, compd. with 1-[2'-(4-methoxyphenyl)[2,5':2',5''-ter-1H-benzimidazol]-5-yl]-3-pyrrolidinamine (1:1) (9CI)  
MF C32 H28 N8 O . Unspecified

CM 1



CM 2

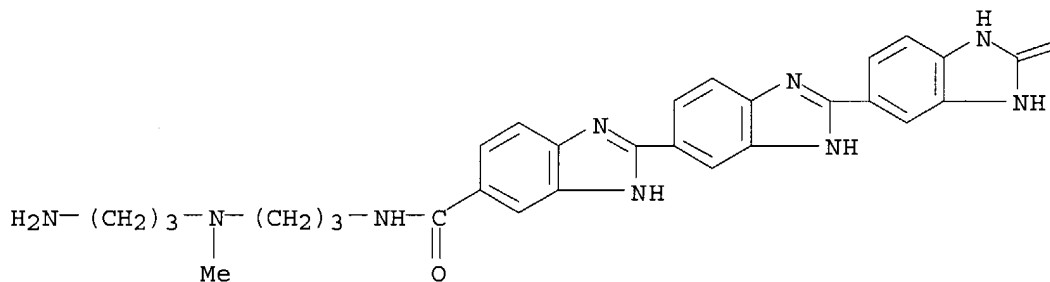
\*\*RELATED SEQUENCES AVAILABLE WITH SEQLINK\*\*

\*\*\* STRUCTURE DIAGRAM IS NOT AVAILABLE \*\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L6 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN [2,5':2',5''-Ter-1H-benzimidazole]-5-carboxamide, N-[3-[(3-aminopropyl)methylamino]propyl]-2'',3''-dihydro-2''-oxo- (9CI)  
MF C29 H31 N9 O2

PAGE 1-A



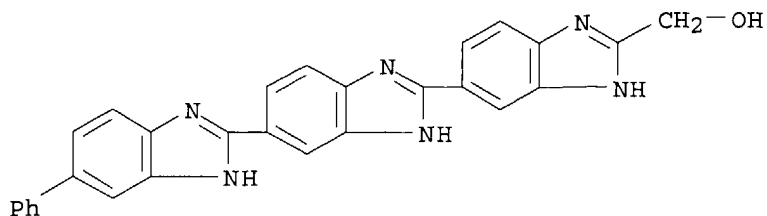
PAGE 1-B

=O

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

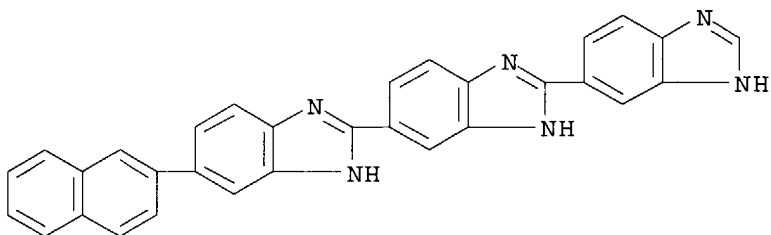
L6 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN [2,5':2',5''-Ter-1H-benzimidazole]-2''-methanol, 5-phenyl- (9CI)  
MF C28 H20 N6 O



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

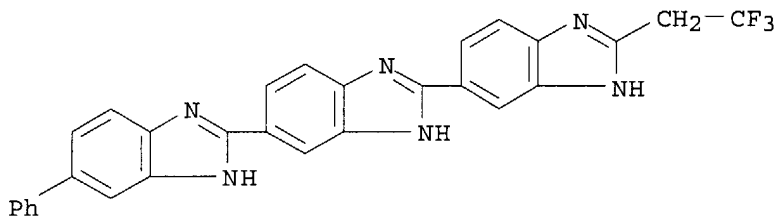
L6 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
 IN 2,5':2'',5'''-Ter-1H-benzimidazole, 5-(2-naphthalenyl)- (9CI)  
 MF C31 H20 N6



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

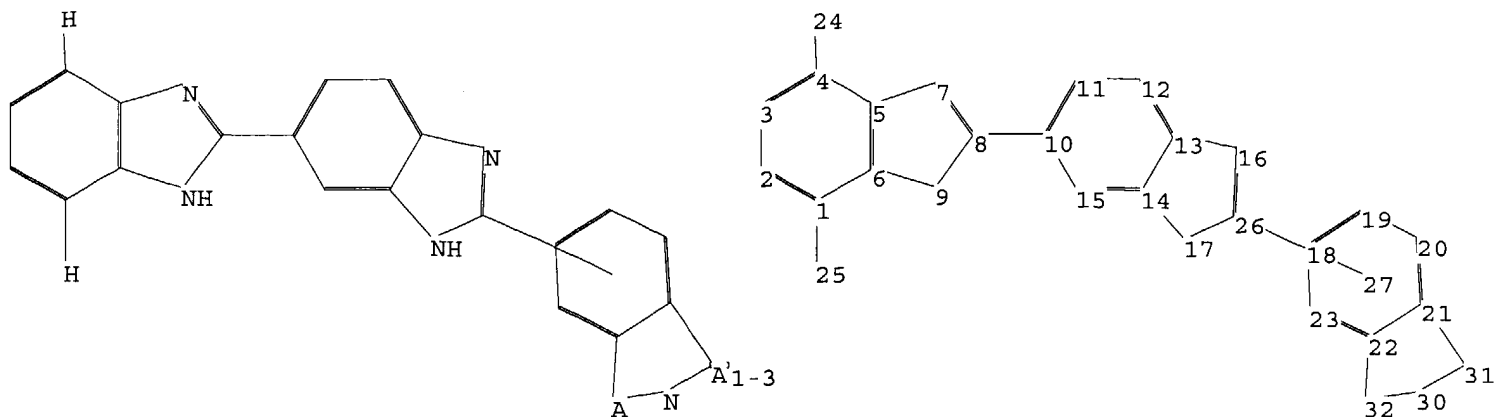
L6 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
 IN 2,5':2'',5'''-Ter-1H-benzimidazole, 5-phenyl-2''-(2,2,2-trifluoroethyl)- (9CI)  
 MF C29 H19 F3 N6



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=>  
 Uploading C:\Program Files\Stnexp\Queries\10690800g.str



chain nodes :

24 25

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 26 30 31 32

chain bonds :

1-25 4-24 8-10

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 10-11 10-15 11-12 12-13 13-14 13-16 14-15 14-17 16-26 17-26 18-23 18-19 19-20 20-21 21-22 21-31 22-23 22-32 30-31 30-32

exact/norm bonds :

5-7 6-9 7-8 8-9 13-16 14-17 16-26 17-26 21-31 22-32 30-31 30-32

exact bonds :

1-25 4-24 8-10

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 10-11 10-15 11-12 12-13 13-14 14-15 18-23 18-19 19-20 20-21 21-22 22-23

G1

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom 22:Atom 23:Atom 24:CLASS 25:CLASS 26:Atom 27:CLASS 30:Atom 31:Atom 32:Atom

L7 STRUCTURE UPLOADED

=> s l7

SAMPLE SEARCH INITIATED 14:06:00 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 54 TO ITERATE

100.0% PROCESSED 54 ITERATIONS

0 ANSWERS

SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 640 TO 1520

PROJECTED ANSWERS: 0 TO 0

L8 0 SEA SSS SAM L7

=> s l7 full

FULL SEARCH INITIATED 14:06:04 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 1067 TO ITERATE

100.0% PROCESSED 1067 ITERATIONS

1 ANSWERS

SEARCH TIME: 00.00.01

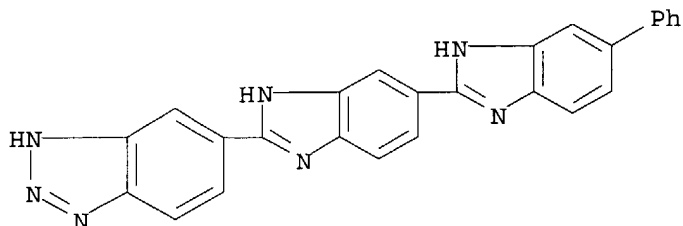
L9 1 SEA SSS FUL L7

=> d scan

L9 1 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

IN 1H-Benzotriazole, 5-(5-phenyl[2,5'-bi-1H-benzimidazol]-2'-yl)- (9CI)

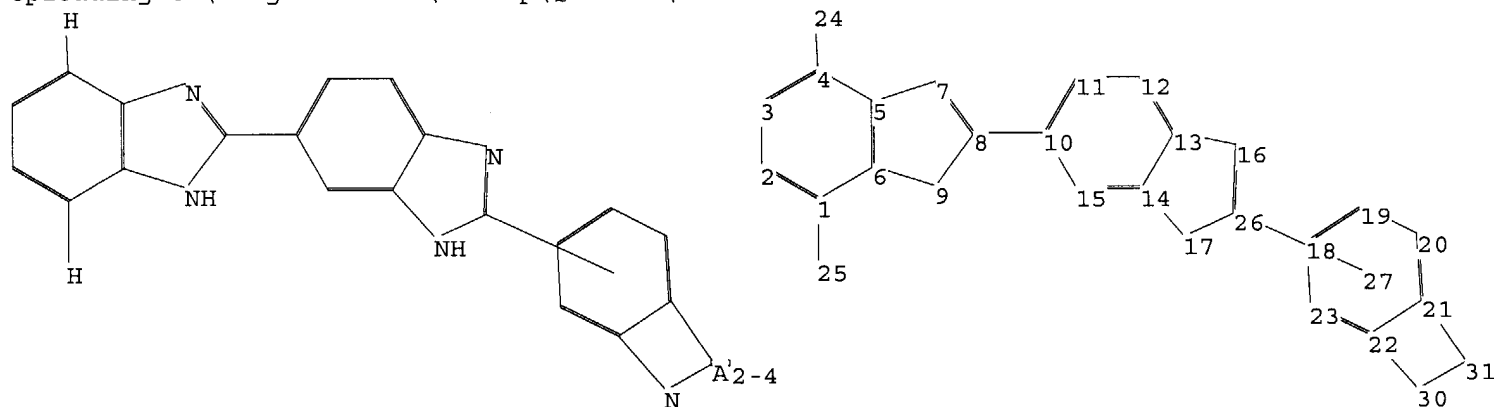
MF C26 H17 N7



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

ALL ANSWERS HAVE BEEN SCANNED

=>  
Uploading C:\Program Files\Stnexp\Queries\10690800h.str



chain nodes :

24 25

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 26 30 31

chain bonds :

1-25 4-24 8-10

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 10-11 10-15 11-12 12-13 13-14 13-16  
14-15 14-17 16-26 17-26 18-23 18-19 19-20 20-21 21-22 21-31 22-23 22-30 30-31

exact/norm bonds :

5-7 6-9 7-8 8-9 13-16 14-17 16-26 17-26 21-31 22-30 30-31

exact bonds :

1-25 4-24 8-10

normalized bonds :

1-2 1-6 2-3 3-4 4-5 5-6 10-11 10-15 11-12 12-13 13-14 14-15 18-23 18-19 19-20  
20-21 21-22 22-23

G1

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom  
12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom  
22:Atom 23:Atom 24:CLASS 25:CLASS 26:Atom 27:CLASS 30:Atom 31:Atom

L10 STRUCTURE UPLOADED

=> s l10

SAMPLE SEARCH INITIATED 14:06:58 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 14 TO ITERATE

100.0% PROCESSED 14 ITERATIONS  
SEARCH TIME: 00.00.01

4 ANSWERS

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 56 TO 504  
PROJECTED ANSWERS: 4 TO 200

L11 4 SEA SSS SAM L10

=> s l10 full

FULL SEARCH INITIATED 14:07:02 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 364 TO ITERATE

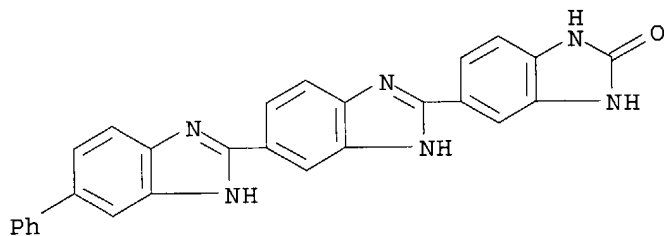
100.0% PROCESSED 364 ITERATIONS  
SEARCH TIME: 00.00.01

75 ANSWERS

L12 75 SEA SSS FUL L10

=> d scan

L12 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN [2,5':2',5''-Ter-1H-benzimidazol]-2''(3''H)-one, 5-phenyl- (9CI)  
MF C27 H18 N6 O

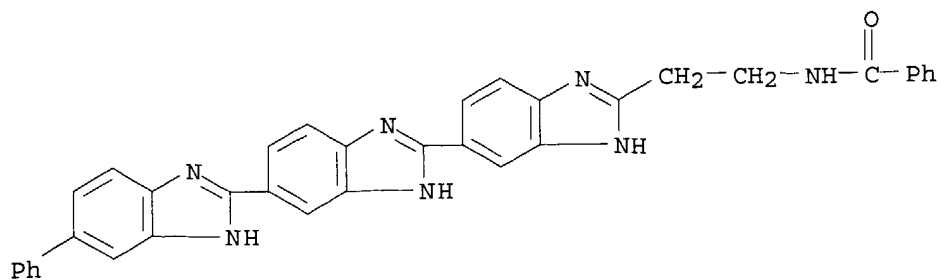


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L12 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN Benzamide, N-[2-(5-phenyl[2,5':2',5''-ter-1H-benzimidazol]-2''-yl)ethyl]-  
(9CI)  
MF C36 H27 N7 O

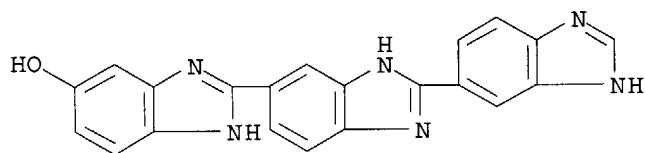




\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

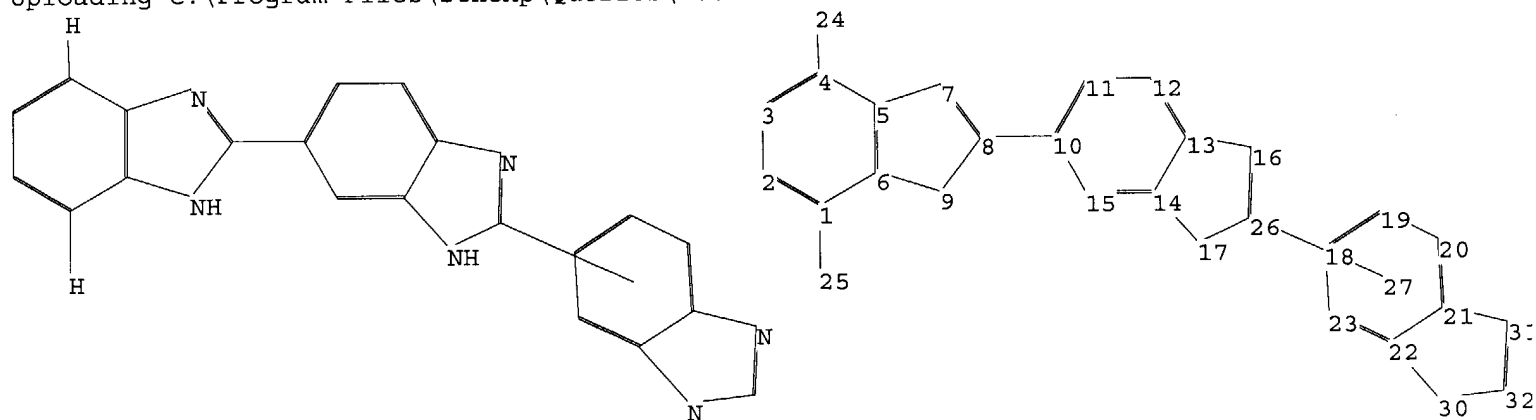
L12 75 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
 IN [2,5':2',5''-Ter-1H-benzimidazol]-5-ol (9CI)  
 MF C21 H14 N6 O



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=>  
 Uploading C:\Program Files\Stnexp\Queries\10690800i.str



chain nodes :

24 25

ring nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 19 20 21 22 23 26 30 31  
 32

chain bonds :

1-25 4-24 8-10

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-9 7-8 8-9 10-11 10-15 11-12 12-13 13-14 13-16  
 14-15 14-17 16-26 17-26 18-23 18-19 19-20 20-21 21-22 21-31 22-23 22-30 30-32  
 31-32

exact/norm bonds :

5-7 6-9 7-8 8-9 13-16 14-17 16-26 17-26 21-31 22-30 30-32 31-32  
exact bonds :  
1-25 4-24 8-10  
normalized bonds :  
1-2 1-6 2-3 3-4 4-5 5-6 10-11 10-15 11-12 12-13 13-14 14-15 18-23 18-19 19-20  
20-21 21-22 22-23

G1

Match level :

1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom 11:Atom  
12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:Atom 20:Atom 21:Atom  
22:Atom 23:Atom 24:CLASS 25:CLASS 26:Atom 27:CLASS 30:Atom 31:Atom 32:Atom

L13 STRUCTURE UPLOADED

=> s l13

SAMPLE SEARCH INITIATED 14:07:31 FILE 'REGISTRY'  
SAMPLE SCREEN SEARCH COMPLETED - 15 TO ITERATE

100.0% PROCESSED 15 ITERATIONS 4 ANSWERS  
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*  
PROJECTED ITERATIONS: 68 TO 532  
PROJECTED ANSWERS: 4 TO 200

L14 4 SEA SSS SAM L13

=> s l13 full

FULL SEARCH INITIATED 14:07:34 FILE 'REGISTRY'  
FULL SCREEN SEARCH COMPLETED - 418 TO ITERATE

100.0% PROCESSED 418 ITERATIONS 70 ANSWERS  
SEARCH TIME: 00.00.01

L15 70 SEA SSS FUL L13

=> d his

(FILE 'HOME' ENTERED AT 14:04:20 ON 21 JUN 2004)

FILE 'REGISTRY' ENTERED AT 14:04:25 ON 21 JUN 2004

L1 STRUCTURE UPLOADED  
L2 0 S L1  
L3 0 S L1 FULL  
L4 STRUCTURE UPLOADED  
L5 4 S L4  
L6 75 S L4 FULL  
L7 STRUCTURE UPLOADED  
L8 0 S L7  
L9 1 S L7 FULL  
L10 STRUCTURE UPLOADED  
L11 4 S L10  
L12 75 S L10 FULL  
L13 STRUCTURE UPLOADED  
L14 4 S L13  
L15 70 S L13 FULL

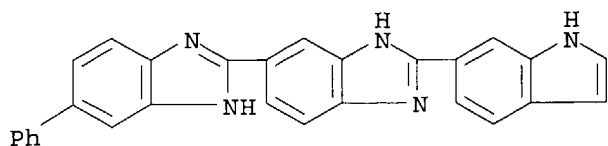
=> s l12 not l15

L16 5 L12 NOT L15

=> d scan

L16 5 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN

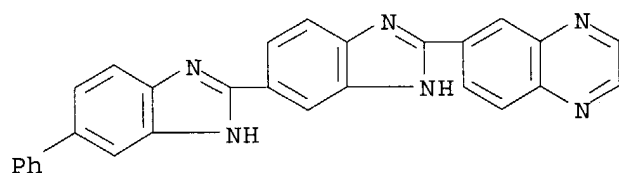
IN 2,5'-Bi-1H-benzimidazole, 2'-(1H-indol-6-yl)-5-phenyl- (9CI)  
MF C28 H19 N5



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

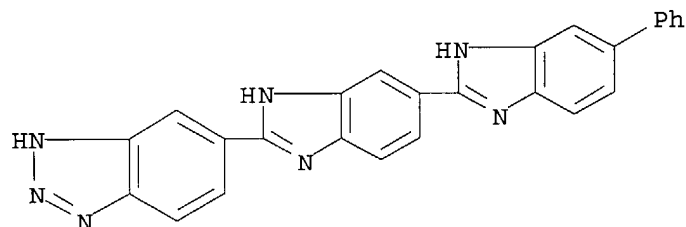
L16 5 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN Quinoxaline, 6-(5-phenyl[2,5'-bi-1H-benzimidazol]-2'-yl)- (9CI)  
MF C28 H18 N6



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

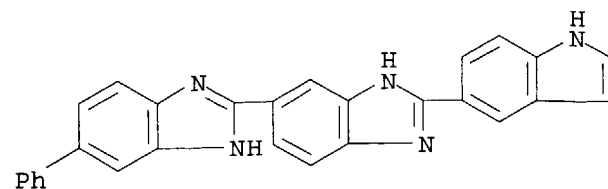
L16 5 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN 1H-Benzotriazole, 5-(5-phenyl[2,5'-bi-1H-benzimidazol]-2'-yl)- (9CI)  
MF C26 H17 N7



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

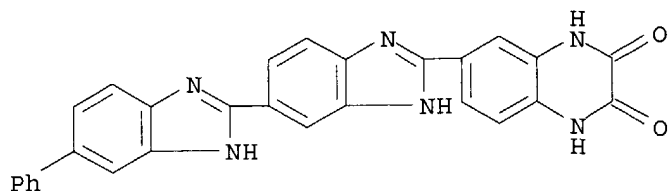
L16 5 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN 2,5'-Bi-1H-benzimidazole, 2'-(1H-indol-5-yl)-5-phenyl- (9CI)  
MF C28 H19 N5



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L16 5 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN  
IN 2,3-Quinoxalinedione, 1,4-dihydro-6-(5-phenyl[2,5'-bi-1H-benzimidazol]-2'-yl)-(9CI)  
MF C28 H18 N6 O2



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

ALL ANSWERS HAVE BEEN SCANNED

=> fil medline caplus  
COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
777.52	777.73

FULL ESTIMATED COST

FILE 'MEDLINE' ENTERED AT 14:08:11 ON 21 JUN 2004

FILE 'CAPLUS' ENTERED AT 14:08:11 ON 21 JUN 2004  
USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.  
PLEASE SEE "HELP USAGETERMS" FOR DETAILS.  
COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

=> d his

(FILE 'HOME' ENTERED AT 14:04:20 ON 21 JUN 2004)

FILE 'REGISTRY' ENTERED AT 14:04:25 ON 21 JUN 2004

L1 STRUCTURE UPLOADED  
L2 0 S L1  
L3 0 S L1 FULL  
L4 STRUCTURE UPLOADED  
L5 4 S L4  
L6 75 S L4 FULL  
L7 STRUCTURE UPLOADED  
L8 0 S L7  
L9 1 S L7 FULL  
L10 STRUCTURE UPLOADED  
L11 4 S L10  
L12 75 S L10 FULL  
L13 STRUCTURE UPLOADED  
L14 4 S L13  
L15 70 S L13 FULL  
L16 5 S L12 NOT L15

FILE 'MEDLINE, CAPLUS' ENTERED AT 14:08:11 ON 21 JUN 2004

=> s 16 or 19 or l16  
L17 28 L6 OR L9 OR L16

=> s 117 and py<1998  
L18 7 L17 AND PY<1998

=> d 1-7

L18 ANSWER 1 OF 7 CAPLUS COPYRIGHT 2004 ACS on STN  
AN 1997:802609 CAPLUS  
DN 128:123523  
TI A terbenzimidazole that preferentially binds and conformationally alters structurally distinct DNA duplex domains: a potential mechanism for topoisomerase I poisoning  
AU Pilch, Daniel S.; Xu, Zhitao; Sun, Qun; LaVoie, Edmond J.; Liu, Leroy F.; Breslauer, Kenneth J.  
CS Dep. Pharmacol., Univ. Med. Dentistry New Jersey, Piscataway, NJ, 08854, USA  
SO Proceedings of the National Academy of Sciences of the United States of America (1997), 94(25), 13565-13570  
CODEN: PNASA6; ISSN: 0027-8424  
PB National Academy of Sciences  
DT Journal  
LA English  
RE.CNT 45 THERE ARE 45 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L18 ANSWER 2 OF 7 CAPLUS COPYRIGHT 2004 ACS on STN  
AN 1997:526711 CAPLUS  
DN 127:117067  
TI Terbenzimidazoles: Influence of 2''-, 4-, and 5-Substituents on Cytotoxicity and Relative Potency as Topoisomerase I Poisons  
AU Kim, Jung Sun; Yu, Chiang; Liu, Angela; Liu, Leroy F.; LaVoie, Edmond J.  
CS Department of Pharmaceutical Chemistry, Rutgers The State University of New Jersey, Piscataway, NJ, 08855, USA  
SO Journal of Medicinal Chemistry (1997), 40(18), 2818-2824  
CODEN: JMCMAR; ISSN: 0022-2623  
PB American Chemical Society  
DT Journal  
LA English

L18 ANSWER 3 OF 7 CAPLUS COPYRIGHT 2004 ACS on STN  
AN 1997:299447 CAPLUS  
DN 126:340359  
TI Differential Poisoning of Human and Aspergillus nidulans DNA Topoisomerase I by Bi- and Terbenzimidazoles  
AU Goldman, Gustavo H.; Yu, Chiang; Wu, Hong-Yan; Sanders, Marilyn M.; LaVoie, Edmond J.; Liu, Leroy F.  
CS Department of Pharmacology Robert Wood Johnson Medical School, University of Medicine and Dentistry of New Jersey, Piscataway, NJ, USA  
SO Biochemistry (1997), 36(21), 6488-6494  
CODEN: BICHAW; ISSN: 0006-2960  
PB American Chemical Society  
DT Journal  
LA English

L18 ANSWER 4 OF 7 CAPLUS COPYRIGHT 2004 ACS on STN  
AN 1997:41984 CAPLUS  
DN 126:59953  
TI Preparation of tribenzimidazoles useful as topoisomerase I inhibitors.  
IN Lavoie, Edmond J.; Liu, Leroy Fong; Sun, Qun  
PA Rutgers, the State University of New Jersey, USA; Lavoie, Edmond J.; Liu, Leroy Fong; Sun, Qun  
SO PCT Int. Appl., 36 pp.  
CODEN: PIXXD2  
DT Patent  
LA English  
FAN.CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 9636612	A1	19961121	WO 1996-US6853	19960514 <--
	W: AL, AM, AT, AU, AZ, BB, BG, BR, BY, CA, CH, CN, CZ, DE, DK, EE, ES, FI, GB, GE, HU, IS, JP, KE, KG, KP, KR, KZ, LK, LR, LS, LT,				

LU, LV, MD, MG, MK, MN, MW, MX, NO, NZ, PL, PT, RO, RU, SD, SE,  
 SG, SI  
 RW: KE, LS, MW, SD, SZ, UG, AT, BE, CH, DE, DK, ES, FI, FR, GB, GR,  
 IE, IT, LU, MC, NL, PT, SE, BF, BJ, CF, CG, CI, CM, GA, GN  
 US 5807874 A 19980915 US 1995-442657 19950517  
 US 5767142 A 19980616 US 1996-618988 19960320  
 CA 2221248 AA 19961121 CA 1996-2221248 19960514 <--  
 AU 9657466 A1 19961129 AU 1996-57466 19960514 <--  
 AU 713317 B2 19991125  
 EP 839140 A1 19980506 EP 1996-915784 19960514  
 R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT,  
 IE, SI, LT, LV, FI  
 JP 11508229 T2 19990721 JP 1996-534375 19960514  
 BR 9608476 A 19990817 BR 1996-8476 19960514  
 US 5948797 A 19990907 US 1997-782064 19970113  
 AU 9952697 A1 19991202 AU 1999-52697 19991005  
 AU 730456 B2 20010308  
 PRAI US 1995-442657 A 19950517  
 US 1996-618988 A 19960320  
 AU 1996-57466 A3 19960514  
 WO 1996-US6853 W 19960514  
 OS MARPAT 126:59953  
  
 L18 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2004 ACS on STN  
 AN 1996:607536 CAPLUS  
 DN 125:265169  
 TI Isohelicity and Phasing in Drug-DNA Sequence Recognition: Crystal  
 Structure of a Tris(benzimidazole)-Oligonucleotide Complex  
 AU Clark, George R.; Gray, Emily J.; Neidle, Stephen; Li, Yu-Hua; Leupin,  
 Werner  
 CS CRC Biomolecular Structure Unit, Institute of Cancer Research,  
 Sutton/Surrey, SM2 5NG, UK  
 SO Biochemistry (1996), 35(43), 13745-13752  
 CODEN: BICHAW; ISSN: 0006-2960  
 PB American Chemical Society  
 DT Journal  
 LA English  
  
 L18 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2004 ACS on STN  
 AN 1996:405834 CAPLUS  
 DN 125:157872  
 TI Synthesis and structure-activity relationships of novel mammalian DNA  
 topoisomerase I inhibitors (Hoechst dyes, terbenzimidazoles, antitumor  
 drugs)  
 AU Sun, Qun  
 CS Rutgers, State Univ., New Brunswick, NJ, USA  
 SO (1996) 173 pp. Avail.: Univ. Microfilms Int., Order No.  
 DA9618556  
 From: Diss. Abstr. Int., B 1996, 57(2), 1093  
 DT Dissertation  
 LA English  
  
 L18 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2004 ACS on STN  
 AN 1995:746950 CAPLUS  
 DN 123:198740  
 TI Synthesis and Evaluation of Terbenzimidazoles as Topoisomerase I  
 Inhibitors  
 AU Sun, Qun; Gatto, Barbara; Yu, Chiang; Liu, Angela; Liu, Leroy F.; LaVoie,  
 Edmond J.  
 CS Department of Pharmaceutical Chemistry, Rutgers, State University of New  
 Jersey, Piscataway, NJ, 08855, USA  
 SO Journal of Medicinal Chemistry (1995), 38(18), 3638-44  
 CODEN: JMCMAR; ISSN: 0022-2623  
 PB American Chemical Society  
 DT Journal  
 LA English  
 OS CASREACT 123:198740